

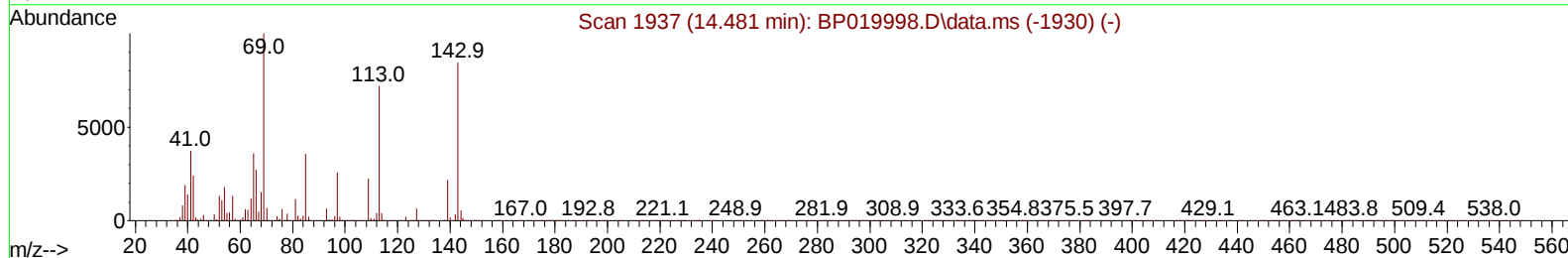
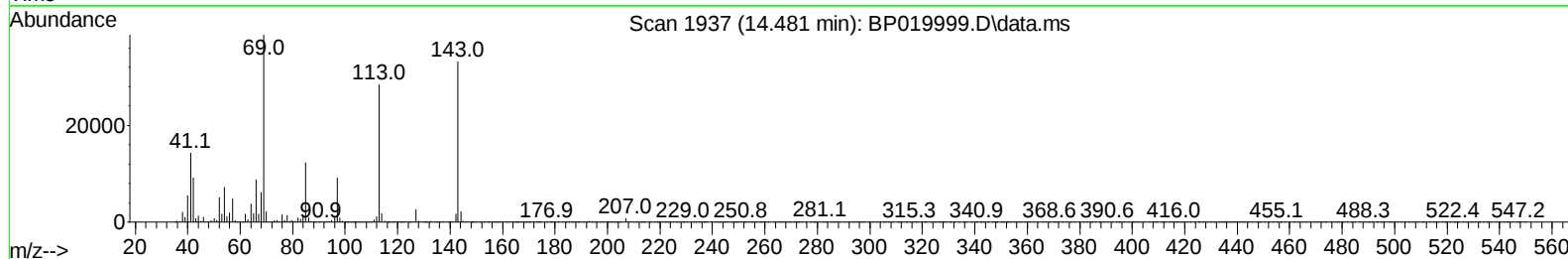
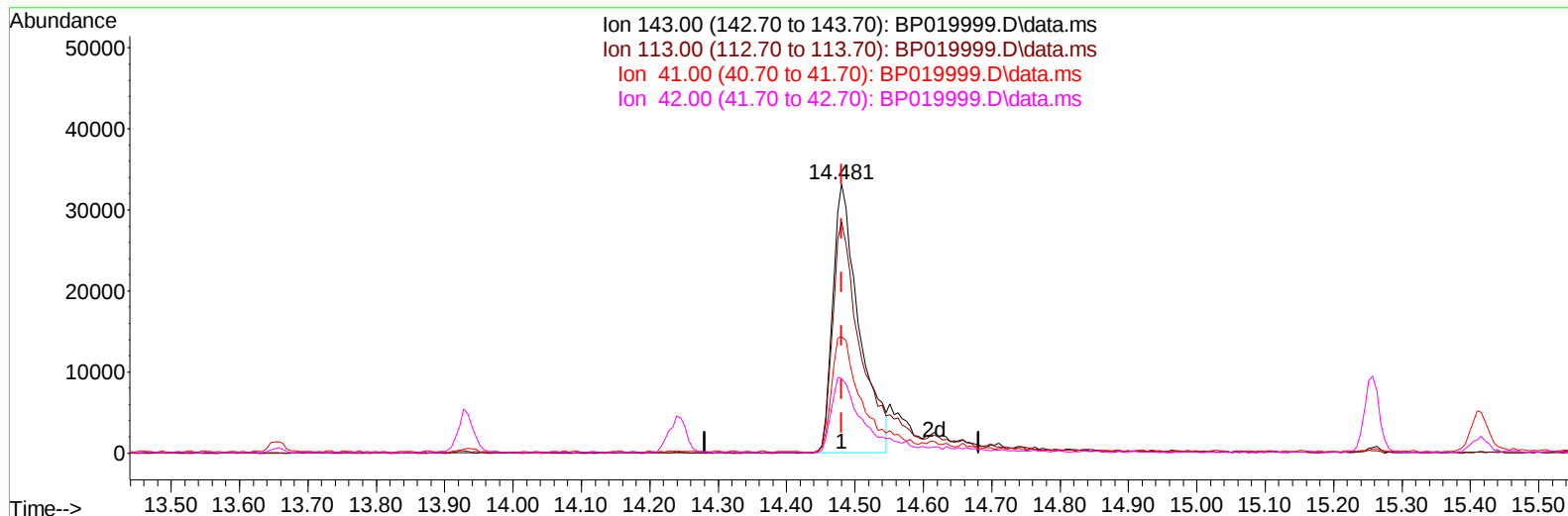
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032624\
Data File : BP019999.D
Acq On : 26 Mar 2024 16:15
Operator : MA/JU
Sample : PB159783BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK783

Manual IntegrationsAPPROVED

Quant Time: Mar 26 17:20:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 25 10:46:59 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
Supervised By :mohammad ahmed 03/29/2024



TIC: BP019999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (-0.000) 31.39 ng/u1

response 88970

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	85.82
41.00	40.90	43.28
42.00	28.50	27.70

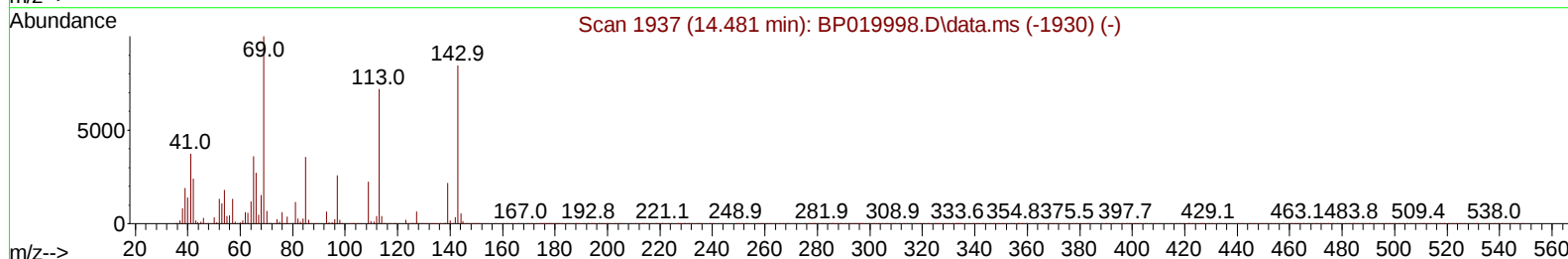
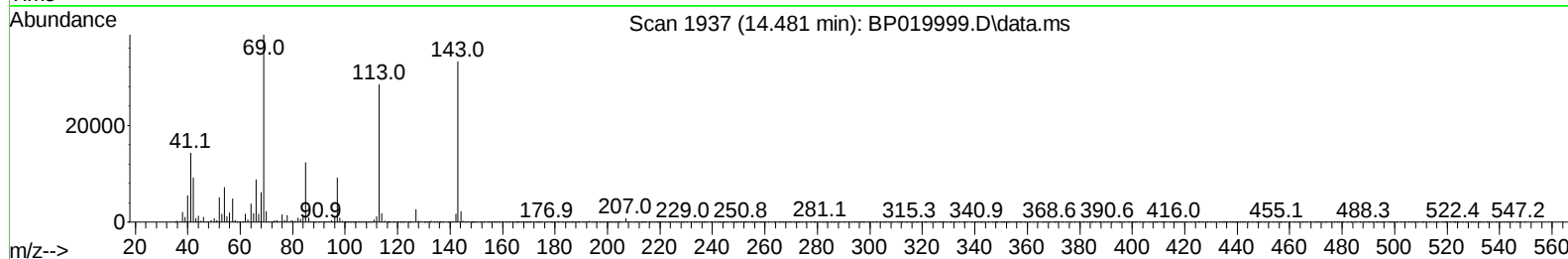
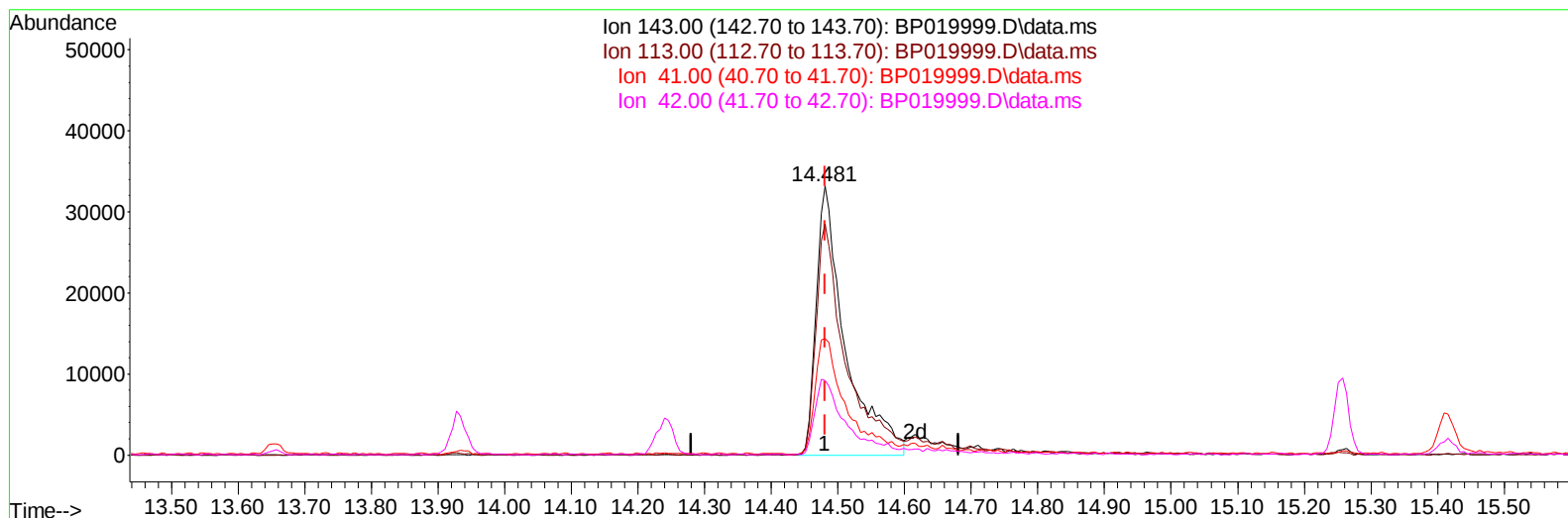
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TIC: BP019999.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (-0.000) 35.58 ng/ul m

response 100820

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	85.82
41.00	40.90	43.28
42.00	28.50	27.70

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :mohammad ahmed 03/29/2024

Quant Time: Mar 27 00:02:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	90793	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.352	136	362551	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.240	164	257110	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	574835	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	567473	20.000	ng/ul	-0.01
88) Perylene-d12	24.827	264	521838	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	14533	7.013	ng/uL	-0.01
4) Pyridine-d5	3.634	84	152591	24.387	ng/ul	0.00
7) Phenol-d5	6.781	99	243956	31.859	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	6.958	67	168660	33.807	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.134	132	180912	32.898	ng/ul	-0.02
15) 4-Methylphenol-d8	8.293	113	184024	31.062	ng/ul	-0.01
21) Nitrobenzene-d5	8.746	128	91589	32.959	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.457	143	102299	32.902	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.993	165	158492	27.912	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.651	166	656981	36.445	ng/ul	0.00
49) Acenaphthylene-d8	13.934	160	665651	31.610	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	100820m	35.576	ng/ul	0.00
60) Fluorene-d10	15.257	176	529831	34.503	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	106899	31.067	ng/ul	0.00
73) Anthracene-d10	17.163	188	904052	35.529	ng/ul	0.00
81) Pyrene-d10	19.563	212	1151510	34.586	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.592	264	873003	34.823	ng/ul	-0.02

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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